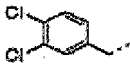
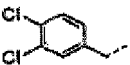
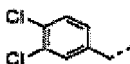
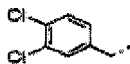
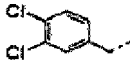
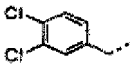
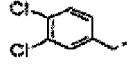

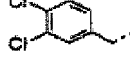
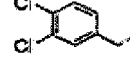



AMENDMENTS TO THE SPECIFICATION

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


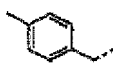
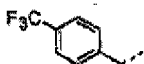
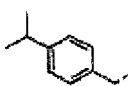
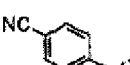
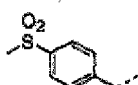
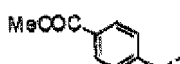
24

《Table 1-1》 X=Single Bond, q=0, r=0, Y=-(R4)C=C(R5)-							
Compound No.	R1-(CH2)p--	R2	R3	R4	R5	R6	R7
1-1-1		H	H	H	H	H	H
1-1-2		H	H	H	H	NO2	H
1-1-3		H	H	H	H	Me	H
1-1-4		H	H	H	H	Cl	H
1-1-5		H	H	H	H	F	H
1-1-6		H	H	H	H	CF3	H
1-1-7		H	H	H	H	COOH	H
1-1-8		H	H	H	Cl	Cl	H
1-1-9		H	H	H	H	H	Me
1-1-10		H	H	H	H	MeO	H
1-1-11		H	H	H	H	H	NO2

Please delete page 37 and replace with the following page 37:


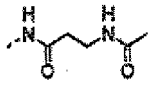
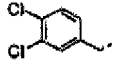
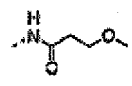

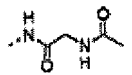
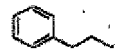
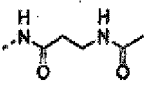

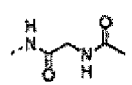
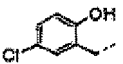
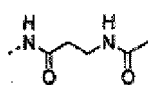

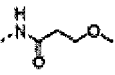
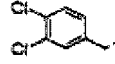
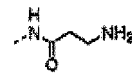
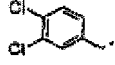
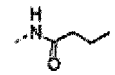
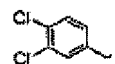
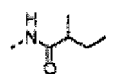
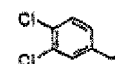
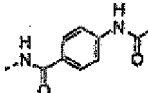
37

《Table 1-2》 X = Single Bond, q = 0, r = 0, Y = -(R4)C(R5)-

Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R6	R7
1-2-1		H	H	H	H	H	H
1-2-2		H	H	H	H	H	H
1-2-3		H	H	H	H	H	H
1-2-4		H	H	H	H	H	H
1-2-5		H	H	H	H	H	H
1-2-6		H	H	H	H	H	H
1-2-7		H	H	H	H	H	H
1-2-8		H	H	H	H	H	H
1-2-9		H	H	H	H	H	H
1-2-10		H	H	H	H	H	H

Please delete page 54 and replace with the following page 54:

54

$\langle \text{Table 1-3} \rangle$ X = Single Bond, q = 0, r = 0, Y = $-(R_4)C=C(R_5)-$						
Compound No.	R1-(CH ₂) _p -	R2	R3	R4	R5	R7
1-3-1		H	H	H	H	
1-3-2		H	H	H	H	
1-3-3		H	H	H	H	
1-3-4		H	H	H	H	
1-3-5		H	H	H	H	
1-3-6		H	H	H	H	
1-3-7		H	H	H	H	
1-3-8		H	H	H	H	
1-3-9		H	H	H	H	
1-3-10		H	H	H	H	
1-3-11		H	H	H	H	

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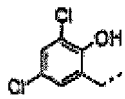
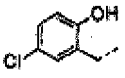


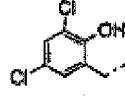
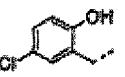



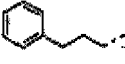
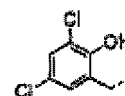
81

(Table 1-4) X = Single Bond, q = 0, r = 0, Y = -(R4)O=C(R5)-

Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R6	R7
1-4-1		H	H	H	H	H	
1-4-2		H	H	H	H	H	
1-4-3		H	H	H	H	H	
1-4-4		H	H	H	H	H	
1-4-5		H	H	H	H	H	
1-4-6		H	H	H	H	H	
1-4-7		H	H	H	H	H	
1-4-8		H	H	H	H	H	
1-4-9		H	H	H	H	H	
1-4-10		H	H	H	H	H	
1-4-11		H	H	H	H	H	

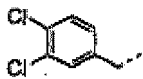

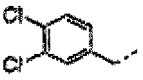
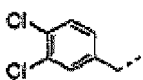
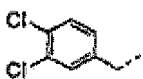





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88

Table 1-5 X = Single Bond, q = 0, r = 0, Y = -(R4)C=C(R5)-							
Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R6	R7
1-5-1		H	H	H	H	COOEt	H
1-5-2		H	H	H	H	COOEt	H
1-5-3		H	H	H	H	COOEt	H
1-5-4		H	H	H	H	COOEt	H
1-5-5		H	H	H	H	COOCH(Me)2	H
1-5-6		H	H	H	H	COOCH(Me)2	H
1-5-7		H	H	H	H	COOCH(Me)2	H
1-5-8		H	H	H	H	COOCH(Me)2	H
1-5-9		H	H	H	H	COOMe	H
1-5-10		H	H	H	H	COOMe	H
1-5-11		H	H	H	H	COOMe	H

Please delete page 107 and replace with the following page 107:

107

Table 1-6 X = Single Bond, $\alpha = 0$, $r = 0$, Y = $-(R_4)C=O(R_5)-$							
Compound No.	R1-(CH ₂) _p -	R2	R3	R4	R5	R6	R7
1-6-1		H	Et	H	H	H	H
1-6-2		Et	Et	H	H	H	H
1-6-3		CH ₂ C ₆ H ₅	CH ₂ C ₆ H ₅	H	H	H	H
1-6-4		H	CH ₂ C ₆ H ₅	H	H	H	H
1-6-5		(CH ₂) ₅ CH ₃	(CH ₂) ₅ CH ₃	H	H	H	H
1-6-6		H	(CH ₂) ₅ CH ₃	H	H	H	H
1-6-7		(CH ₂) ₃ C ₆ H ₅	(CH ₂) ₃ C ₆ H ₅	H	H	H	H
1-6-8		H	(CH ₂) ₃ C ₆ H ₅	H	H	H	H
1-6-9		H	CH ₂ COOMe	H	H	H	H
1-6-10		H	(CH ₂) ₄ COOEt	H	H	H	H